

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031119\
 Data File : BE098874.D
 Acq On : 11 Mar 2019 23:51
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 12 16:38:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 2019
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	112	0.00
2	1,4-Dioxane	0.400	0.375	6.3	109	0.00
3	n-Nitrosodimethylamine	0.400	0.240	40.0#	66	0.03
4 S	2-Fluorophenol	0.400	0.356	11.0	97	0.00
5 S	Phenol-d6	0.400	0.399	0.3	112	0.00
6	bis(2-Chloroethyl)ether	0.400	0.324	19.0	95	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	125	0.00
8 S	Nitrobenzene-d5	0.400	0.368	8.0	128	0.00
9	Naphthalene	0.400	0.400	0.0	128	0.00
10	Hexachlorobutadiene	0.400	0.371	7.3	120	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.397	0.8	127	0.00
12	2-Methylnaphthalene	0.400	0.383	4.3	124	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	132	0.00
14 S	2,4,6-Tribromophenol	0.400	0.343	14.2	114	0.00
15 S	2-Fluorobiphenyl	0.400	0.406	-1.5	136	0.00
16	Acenaphthylene	0.400	0.392	2.0	136	0.00
17	Acenaphthene	0.400	0.396	1.0	131	-0.01
18	Fluorene	0.400	0.379	5.3	128	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	127	0.00
20	4-Bromophenyl-phenylether	0.400	0.352	12.0	114	0.00
21	Hexachlorobenzene	0.400	0.382	4.5	128	0.00
22	Pentachlorophenol	0.400	0.434	-8.5	123	0.00
23	Phenanthrene	0.400	0.380	5.0	124	0.00
24	Anthracene	0.400	0.369	7.8	124	0.00
25 SURR	Fluoranthene-d10	0.400	0.357	10.8	118	0.00
26	Fluoranthene	0.400	0.357	10.8	118	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	116	-0.01
28	Pvrene	0.400	0.404	-1.0	117	0.00
29 S	Terphenyl-d14	0.400	0.394	1.5	115	0.00
30	Benzo(a)anthracene	0.400	0.370	7.5	109	0.00
31	Chrysene	0.400	0.403	-0.8	120	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.381	4.8	120	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.398	0.5	117	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	122	-0.01
35	Benzo(b)fluoranthene	0.400	0.387	3.3	122	0.00
36	Benzo(k)fluoranthene	0.400	0.384	4.0	123	-0.01
37 C	Benzo(a)pyrene	0.400	0.394	1.5	124	0.00
38	Dibenzo(a,h)anthracene	0.400	0.383	4.3	118	-0.01
39	Benzo(a,h,i)perylene	0.400	0.383	4.3	119	-0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031119\
Data File : BE098874.D
Acq On : 11 Mar 2019 23:51
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 12 16:38:35 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 06 15:33:24 2019
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

Compound	Amount	Calc.	%Dev Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0			